

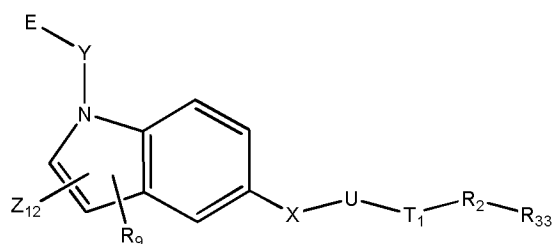
**Amendments to the Claims**

Please amend Claims 88, 92 and 140. The Claim Listing below will replace all prior versions of the claims in the application:

**Claim Listing**

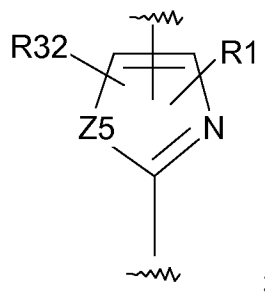
1-48. (Cancelled)

49. (Previously Presented) A compound represented by the following Structural Formula:



or a stereoisomer or a pharmaceutically acceptable salt, wherein:

(a) T1 is selected from the group consisting of



- (b) R1 is selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>8</sub> alkyl, C<sub>1</sub>-C<sub>8</sub> alkenyl, aryl-C<sub>0-4</sub>-alkyl, aryl-C<sub>1-6</sub>-heteroalkyl, heteroaryl-C<sub>0-4</sub>-alkyl, and C<sub>3</sub>-C<sub>6</sub> cycloalkylaryl-C<sub>0-2</sub>-alkyl, wherein C<sub>1</sub>-C<sub>8</sub> alkyl, C<sub>1</sub>-C<sub>8</sub> alkenyl, aryl-C<sub>0-4</sub>-alkyl, aryl-C<sub>1-6</sub>-heteroalkyl, heteroaryl-C<sub>0-4</sub>-alkyl, and C<sub>3</sub>-C<sub>6</sub> cycloalkylaryl-C<sub>0-2</sub>-alkyl are each optionally substituted with from one to three substituents independently selected from R1';
- (c) R1', R26, R27, R28, R31, Z14', and Z15' are each independently selected from the group consisting of hydrogen, hydroxy, cyano, nitro, halo, oxo, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>1</sub>-C<sub>6</sub> alkyl-COOR12, C<sub>1</sub>-C<sub>6</sub> alkoxy, C<sub>1</sub>-C<sub>6</sub> haloalkyl, C<sub>1</sub>-C<sub>6</sub> haloalkyloxy, C<sub>3</sub>-C<sub>7</sub>

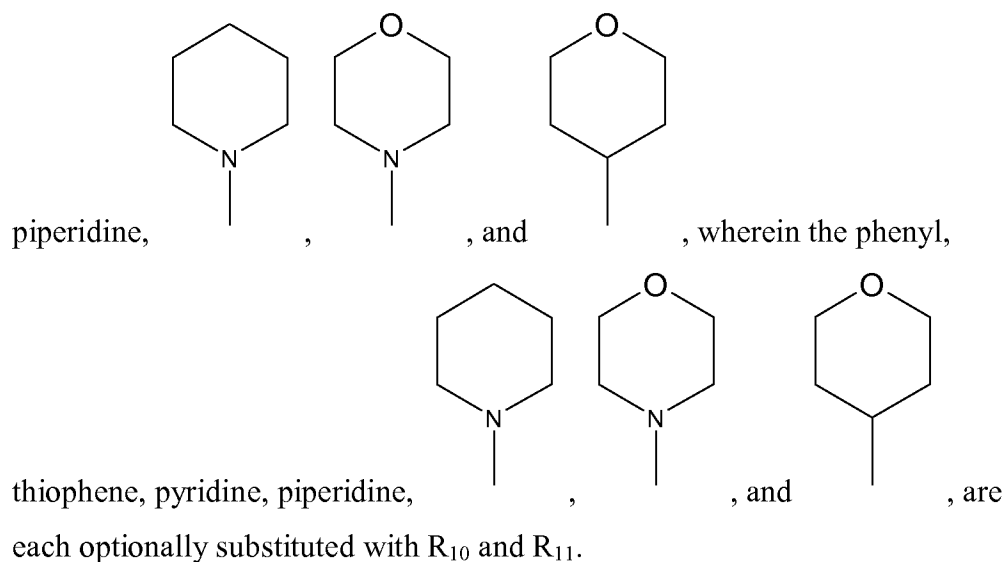
cycloalkyl, aryloxy, aryl-C<sub>0-4</sub>-alkyl, heteroaryl, heterocycloalkyl, C(O)R<sub>13</sub>, COOR<sub>14</sub>, OC(O)R<sub>15</sub>, OS(O)<sub>2</sub>R<sub>16</sub>, N(R<sub>17</sub>)<sub>2</sub>, NR<sub>18</sub>C(O)R<sub>19</sub>, NR<sub>20</sub>SO<sub>2</sub>R<sub>21</sub>, SR<sub>22</sub>, S(O)R<sub>23</sub>, S(O)<sub>2</sub>R<sub>24</sub>, and S(O)<sub>2</sub>N(R<sub>25</sub>)<sub>2</sub>; R<sub>12</sub>, R<sub>13</sub>, R<sub>14</sub>, R<sub>15</sub>, R<sub>16</sub>, R<sub>17</sub>, R<sub>18</sub>, R<sub>19</sub>, R<sub>20</sub>, R<sub>21</sub>, R<sub>22</sub>, R<sub>23</sub>, R<sub>24</sub> and R<sub>25</sub> are each independently selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl and aryl;

- (d) R<sub>2</sub> is selected from the group consisting of C<sub>0</sub>-C<sub>8</sub> alkyl and C<sub>1-6</sub>-heteroalkyl;
- (e) X is O;
- (f) U is -CH<sub>2</sub>-;
- (g) Y is selected from the group consisting of C, O, S, NH and a single bond;
- (h) E is C(R<sub>3</sub>)(R<sub>4</sub>)A and wherein
  - (i) A is selected from the group consisting of carboxyl, tetrazole, C<sub>1</sub>-C<sub>6</sub> alkyl nitrile, carboxamide, sulfonamide and acylsulfonamide; wherein sulfonamide, acylsulfonamide and tetrazole are each optionally substituted with from one to two groups independently selected from R<sup>7</sup>;
  - (ii) each R<sup>7</sup> is independently selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>6</sub> haloalkyl, aryl-C<sub>0</sub>-C<sub>4</sub> alkyl and C<sub>1</sub>-C<sub>6</sub> alkyl, wherein such alkyl and arylalkyl are each optionally substituted with from one to two groups independently selected from R<sup>7'</sup>; each R<sup>7'</sup> is independently selected from halo, C<sub>1</sub>-C<sub>6</sub> alkyl, and haloC<sub>1</sub>-C<sub>6</sub> alkyl;
  - (iii) R<sub>3</sub> is selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>5</sub> alkyl, and C<sub>1</sub>-C<sub>5</sub> alkoxy; and
  - (iv) R<sub>4</sub> is selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>5</sub> alkyl, C<sub>1</sub>-C<sub>5</sub> alkoxy, aryloxy, C<sub>3</sub>-C<sub>6</sub> cycloalkyl, and aryl C<sub>0</sub>-C<sub>4</sub> alkyl, and R<sub>3</sub> and R<sub>4</sub> are optionally combined to form a C<sub>3</sub>-C<sub>4</sub> cycloalkyl, and wherein alkyl, alkoxy, cycloalkyl and aryl-alkyl are each optionally substituted with one to three each independently selected from R<sub>26</sub>;
- (i) Z<sub>5</sub> is O;
- (j) Z<sub>12</sub> is selected from the group consisting of hydrogen and -Z<sub>13</sub>C<sub>0</sub>-C<sub>3</sub>alkylZ<sub>14</sub>;

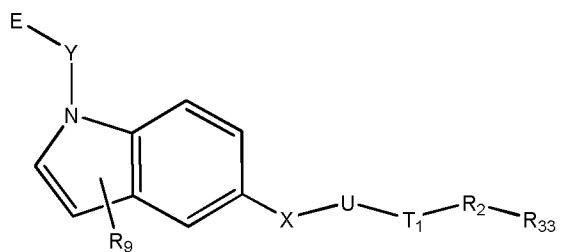
- (k) Z<sub>13</sub> is selected from the group consisting of a single bond, CO, CO<sub>2</sub>, CONZ<sub>15</sub>, and SO<sub>2</sub>;
- (l) Z<sub>14</sub> is selected from the group consisting of aryl and heteroaryl, wherein the aryl and heteroaryl is each optionally substituted with from one to three substituents independently selected from Z<sub>14</sub>';
- (m) Z<sub>15</sub> is selected from the group consisting of hydrogen aryl and heteroaryl, wherein the aryl and heteroaryl is each optionally substituted with from one to three substituents independently selected from Z<sub>15</sub>';
- (n) R<sub>9</sub> is selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>1</sub>-C<sub>4</sub> alkylenyl, halo, aryl-C<sub>0</sub>-C<sub>4</sub> alkyl, heteroaryl, C<sub>1</sub>-C<sub>6</sub> allyl, and OR<sub>29</sub>, and wherein aryl-C<sub>0</sub>-C<sub>4</sub> alkyl, heteroaryl are each optionally substituted with from one to three independently selected from R<sub>27</sub>; R<sub>29</sub> is selected from the group consisting of hydrogen and C<sub>1</sub>-C<sub>4</sub> alkyl;
- (o) R<sub>10</sub>, R<sub>11</sub> are each independently selected from the group consisting of hydrogen, hydroxy, cyano, nitro, halo, oxo, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>0</sub>-C<sub>6</sub> alkyl-COOR<sub>12</sub>'', C<sub>1</sub>-C<sub>6</sub> alkoxy, C<sub>1</sub>-C<sub>6</sub> haloalkyl, C<sub>1</sub>-C<sub>6</sub> haloalkyloxy, C<sub>3</sub>-C<sub>7</sub> cycloalkyl, aryl-C<sub>0-4</sub>-alkyl, aryl-C<sub>1-6</sub>-heteroalkyl, heteroaryl-C<sub>0-4</sub>-alkyl, C<sub>3</sub>-C<sub>6</sub> cycloalkylaryl-C<sub>0-2</sub>-alkyl, aryloxy, C(O)R<sub>13</sub>', COOR<sub>14</sub>', OC(O)R<sub>15</sub>', OS(O)<sub>2</sub>R<sub>16</sub>', N(R<sub>17</sub>')<sub>2</sub>, NR<sub>18</sub>'C(O)R<sub>19</sub>', NR<sub>20</sub>'SO<sub>2</sub>R<sub>21</sub>', SR<sub>22</sub>', S(O)R<sub>23</sub>', S(O)<sub>2</sub>R<sub>24</sub>', and S(O)<sub>2</sub>N(R<sub>25</sub>')<sub>2</sub>; and wherein aryl-C<sub>0-4</sub>-alkyl, aryl- C<sub>1-6</sub>-heteroalkyl, heteroaryl-C<sub>0-4</sub>-alkyl, and C<sub>3</sub>-C<sub>6</sub> cycloalkylaryl-C<sub>0-2</sub>-alkyl are each optionally substituted with from one to three independently selected from R<sub>28</sub>;
- (p) R<sub>12</sub>', R<sub>12</sub>'', R<sub>13</sub>', R<sub>14</sub>', R<sub>15</sub>', R<sub>16</sub>', R<sub>17</sub>', R<sub>18</sub>', R<sub>19</sub>', R<sub>20</sub>', R<sub>21</sub>', R<sub>22</sub>', R<sub>23</sub>', R<sub>24</sub>', and R<sub>25</sub>' are each independently selected from the group consisting of hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl and aryl;
- (q) R<sub>30</sub> is selected from the group consisting of C<sub>1</sub>-C<sub>6</sub> alkyl, aryl-C<sub>0-4</sub>-alkyl, aryl-C<sub>1-6</sub>-heteroalkyl, heteroaryl-C<sub>0-4</sub>-alkyl, and C<sub>3</sub>-C<sub>6</sub> cycloalkylaryl-C<sub>0-2</sub>-alkyl, and wherein C<sub>1</sub>-C<sub>6</sub> alkyl, aryl-C<sub>0-4</sub>-alkyl, aryl- C<sub>1-6</sub>-heteroalkyl, heteroaryl-

C<sub>0-4</sub>-alkyl, and C<sub>3-6</sub> cycloalkylaryl-C<sub>0-2</sub>-alkyl are each optionally substituted with from one to three substituents each independently selected from R<sub>31</sub>;

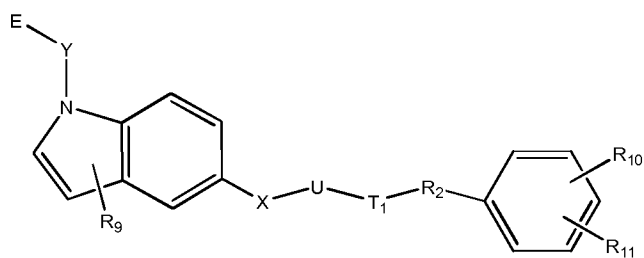
- (r) R<sub>32</sub> is selected from the group consisting of a hydrogen, halo, C<sub>1-6</sub> alkyl, C<sub>1-6</sub> haloalkyl, and C<sub>1-6</sub> alkyloxo;
- (s) R<sub>33</sub> is selected from the group consisting of phenyl, thiophene, pyridine,



50. (Previously Presented) The compound of Claim 49, wherein the compound is represented by the following Structural Formula:

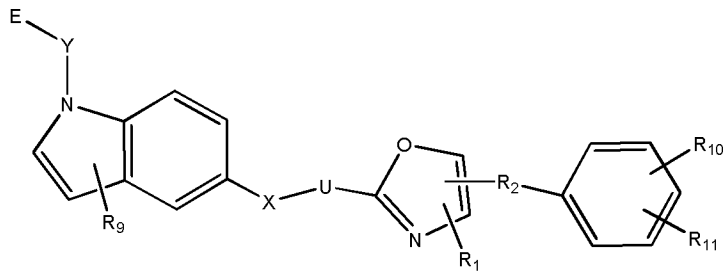


51. (Previously Presented) The compound of Claim 50, wherein the compound is represented by the following Structural Formula:



52-63. (Cancelled)

64. (Previously Presented) The compound of Claim 51, wherein the compound is represented by the following Structural Formula:



65-66. (Cancelled)

67. (Previously Presented) The compound of Claim 64, wherein:

E is C(R3)(R4)-COOH, C<sub>1</sub>-C<sub>6</sub> alkylcarboxyl, or C(R3)(R4)-C<sub>1</sub>-C<sub>6</sub> alkyl-COOH;

R<sub>10</sub> and R<sub>11</sub> are each independently selected from the group consisting of

hydrogen, halo, oxo, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>1</sub>-C<sub>6</sub> alkyl-COOR<sub>12</sub>, C<sub>1</sub>-C<sub>6</sub> alkoxy,

C<sub>1</sub>-C<sub>6</sub> haloalkyl, and C<sub>1</sub>-C<sub>6</sub> haloalkyloxy;

R<sub>9</sub> is selected from the group consisting of hydrogen and C<sub>1</sub>-C<sub>3</sub> alkyl;

R<sub>1</sub>, R<sub>3</sub>, and R<sub>4</sub> are each independently selected from the group consisting of

hydrogen and C<sub>1</sub>-C<sub>2</sub> alkyl; and

R<sub>2</sub> is a bond.

68-87. (Cancelled)

88. (Currently Amended) [[The]] A compound of Claim 49 wherein the compound is selected from the group consisting of:
- {5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-acetic acid;
  - [5-(5-Methyl-2-phenyl-oxazol-4-ylmethoxy)-indol-1-yl]-acetic acid;
  - {5-[2-(4-Fluoro-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-acetic acid;
  - {5-[2-(4-Benzoyloxy-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-acetic acid;
  - 2-Methyl-2-(5-{2-[2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-ethoxy}-indol-1-yl)-propionic acid;
  - {5-[4-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-acetic acid;
  - 2-Methyl-2-{5-[4-methyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-propionic acid;
  - Racemic 2-{5-[4-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-propionic acid;
  - 5-{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-pentanoic acid;
  - 5-{5-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-pentanoic acid;
  - 3-{5-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-propionic acid;
  - {5-[4-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-acetic acid;
  - (5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;
  - (5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;
  - (5-{2-[5-Methyl-2-(tetrahydro-pyran-4-yl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;
  - {5-[2-(2-Butoxy-5-methyl-oxazol-4-yl)-ethoxy]-indol-1-yl}-acetic acid;
  - (5-{2-[2-(5-Bromo-thiophen-2-yl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;
  - {5-[3-(4-Butyl-phenoxy)-propoxy]-indol-1-yl}-acetic acid;
  - (5-{2-[2-(3-Bromo-phenyl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;
  - Racemic 2-(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-propionic acid;
  - (5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

2-(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-2-methyl-propionic acid;

3-(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-propionic acid;

(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-propoxy}-indol-1-yl)-acetic acid;

(S)-(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-propoxy}-indol-1-yl)-acetic acid;

(R)-(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-propoxy}-indol-1-yl)-acetic acid;

Racemic-(5-{1-[4-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-yl]-ethoxy}-indol-1-yl)-acetic acid;

2-Methyl-2-[5-(5-methyl-2-phenyl-oxazol-4-ylmethoxy)-indol-1-yl]-propionic acid;

2-{5-[2-(4-Trifluoromethyl-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-2-methyl-propionic acid;

2-{5-[2-(4-Fluoro-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-2-methyl-propionic acid;

2-{5-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-2-methyl-propionic acid;

2-Methyl-2-(5-{2-[5-methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-propionic acid;

2-(5-{2-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-2-methyl-propionic acid;

2-(5-{2-[2-(5-Bromo-thiophen-2-yl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-2-methyl-propionic acid;

N-(2-{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-acetyl)-methanesulfonamide; and

N-(2-{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-acetyl)-benzenesulfonamide,

or a stereoisomer or a pharmaceutically acceptable salt thereof.

89-91. (Cancelled)

92. (Currently Amended) A method of treating a mammal in need of treatment for a disease selected from the group consisting of diabetes mellitus, Syndrome X, and atherosclerosis, comprising the step of administering to the mammal in need thereof a therapeutically effective amount of the compound of Claim 49, Claim 88 or Claim 140.

93-94. (Cancelled)

95. (Previously Presented) The method of Claim 92, wherein the disease is diabetes mellitus.

96. (Previously Presented) The method of Claim 92, wherein the disease is Syndrome X.

97-139. (Cancelled)

140. (Currently Amended) ~~[[The]]~~ A compound ~~of Claim 49, wherein the compound is~~ selected from the group consisting of:

{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-acetic acid;

[5-(5-Methyl-2-phenyl-oxazol-4-ylmethoxy)-indol-1-yl]-acetic acid;

{5-[2-(4-Fluoro-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-acetic acid;

{5-[2-(4-Benzyloxy-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-acetic acid;

2-Methyl-2-(5-{2-[2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-ethoxy}-indol-1-yl)-propionic acid;

{5-[4-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-acetic acid;

2-Methyl-2-{5-[4-methyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-propionic acid;

Racemic 2-{5-[4-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-propionic acid;



5-{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-pentanoic acid;

5-{5-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-pentanoic acid;

3-{5-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-propionic acid;

{5-[4-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-ylmethoxy]-indol-1-yl}-acetic acid;

(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

(5-{2-[5-Methyl-2-(tetrahydro-pyran-4-yl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

{5-[2-(2-Butoxy-5-methyl-oxazol-4-yl)-ethoxy]-indol-1-yl}-acetic acid;

(5-{2-[2-(5-Bromo-thiophen-2-yl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

(5-{2-[2-(3-Bromo-phenyl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

Racemic 2-(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-propionic acid;

(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-acetic acid;

2-(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-2-methyl-propionic acid;

3-(5-{2-[5-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-propionic acid;

(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-propoxy}-indol-1-yl)-acetic acid;

(S)-(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-propoxy}-indol-1-yl)-acetic acid;

(R)-(5-{2-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-propoxy}-indol-1-yl)-acetic acid;

Racemic-(5-{1-[4-Ethyl-2-(4-trifluoromethyl-phenyl)-oxazol-5-yl]-ethoxy}-indol-1-yl)-acetic acid;

2-Methyl-2-[5-(5-methyl-2-phenyl-oxazol-4-ylmethoxy)-indol-1-yl]-propionic acid;

2-{5-[2-(4-Trifluoromethyl-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-2-methyl-propionic acid;

2-{5-[2-(4-Fluoro-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-2-methyl-propionic acid;

2-{5-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-ylmethoxy]-indol-1-yl}-2-methyl-propionic acid;

2-Methyl-2-(5-{2-[5-methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-yl]-ethoxy}-indol-1-yl)-propionic acid;

2-(5-{2-[2-(4-Bromo-phenyl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-2-methyl-propionic acid;

2-(5-{2-[2-(5-Bromo-thiophen-2-yl)-5-methyl-oxazol-4-yl]-ethoxy}-indol-1-yl)-2-methyl-propionic acid;

N-(2-{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-acetyl)-methanesulfonamide; and

N-(2-{5-[5-Methyl-2-(4-trifluoromethyl-phenyl)-oxazol-4-ylmethoxy]-indol-1-yl}-acetyl)-benzenesulfonamide,<sub>1</sub>

or a stereoisomer or a pharmaceutically acceptable salt thereof.